Ethane, hexafluoro-

Other names:	1,1,1,2,2,2-Hexafluoroethane
	C2F6
	Ethane, 1,1,1,2,2,2-hexafluoro-
	F-116
	FREON 116
	Fluorocarbon 116
	Freon 1166SY
	HFC 116
	Hexafluoroethane
	PERFLUOROETHANE
	R 116
	R-116
	REFRIGERANT-116
	UN 2193
Inchi:	InChI=1S/C2F6/c3-1(4,5)2(6,7)8
InchiKey:	WMIYKQLTONQJES-UHFFFAOYSA-N
Formula:	C2F6
SMILES:	FC(F)(F)C(F)(F)F
Mol. weight [g/mol]:	138.01
CAS:	76-16-4

Physical Properties

Property code	Value	Unit	Source
dm	0.00	debye	KDB
gf	-1258.00	kJ/mol	KDB
hf	-1344.00 ± 5.00	kJ/mol	NIST Webbook
hf	-1330.00	kJ/mol	NIST Webbook
hf	-1331.00 ± 8.00	kJ/mol	NIST Webbook
hf	-1344.00 ± 4.00	kJ/mol	NIST Webbook
hf	-1344.00	kJ/mol	KDB
hfus	4.59	kJ/mol	Joback Method
hvap	12.55	kJ/mol	Joback Method
ie	14.60	eV	NIST Webbook
ie	14.40	eV	NIST Webbook
ie	13.60	eV	NIST Webbook
log10ws	-1.98		Crippen Method
logp	2.111		Crippen Method

mcvol	49.660	ml/mol	McGowan Method
pc	3013.01 ± 3.44	kPa	NIST Webbook
pc	3042.00 ± 8.00	kPa	NIST Webbook
pc	3060.00	kPa	KDB
pc	3040.00 ± 1.10	kPa	NIST Webbook
pt	26.25 ± 0.00	kPa	NIST Webbook
pt	30.26 ± 0.01	kPa	NIST Webbook
rhoc	579.65 ± 19.32	kg/m3	NIST Webbook
rhoc	622.43 ± 9.94	kg/m3	NIST Webbook
rinpol	156.00		NIST Webbook
sl	250.54	J/mol×K	NIST Webbook
tb	195.00	K	KDB
tc	293.03 ± 0.01	K	NIST Webbook
tc	292.80	K	NIST Webbook
tc	292.66 ± 0.05	K	NIST Webbook
tc	293.01 ± 0.01	K	NIST Webbook
tc	292.85 ± 0.05	K	NIST Webbook
tc	293.00	K	KDB
tf	172.40	K	KDB
tf	166.85 ± 0.02	K	NIST Webbook
tf	172.65 ± 0.50	K	NIST Webbook
tf	173.10 ± 0.50	K	NIST Webbook
tt	173.08 ± 0.01	K	NIST Webbook
tt	173.10 ± 0.02	K	NIST Webbook
vc	0.222	m3/kmol	KDB
zc	0.2788500		KDB

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	83.14	J/mol×K	234.32	Joback Method
cpg	114.61	J/mol×K	357.85	Joback Method
cpg	105.32	J/mol×K	316.67	Joback Method
cpg	100.24	J/mol×K	296.08	Joback Method
cpg	94.86	J/mol×K	275.50	Joback Method
cpg	110.11	J/mol×K	337.26	Joback Method
cpg	89.16	J/mol×K	254.91	Joback Method
cpl	131.30	J/mol×K	195.00	NIST Webbook
hfust	2.69	kJ/mol	173.10	NIST Webbook
hfust	3.74	kJ/mol	104.00	NIST Webbook
hfust	2.69	kJ/mol	173.10	NIST Webbook

hsubt	26.00	kJ/mol	103.00	NIST Webbook
hvapt	16.15	kJ/mol	194.87	NIST Webbook
hvapt	16.15	kJ/mol	194.90	KDB
hvapt	17.10	kJ/mol	188.00	NIST Webbook
hvapt	17.30	kJ/mol	186.00	NIST Webbook
hvapt	16.15	kJ/mol	195.00	NIST Webbook
pvap	2270.00	kPa	281.23	Vapor liquid equilibrium data for the hexafluoroethane + carbon dioxide system at temperatures from 253 to 297K and pressures up to 6.5MPa
pvap	1847.00	kPa	273.27	Vapor liquid equilibrium data for the hexafluoroethane + carbon dioxide system at temperatures from 253 to 297K and pressures up to 6.5MPa
pvap	2050.00	kPa	277.24	Vapor liquid equilibrium data for the hexafluoroethane + carbon dioxide system at temperatures from 253 to 297K and pressures up to 6.5MPa
pvap	1659.00	kPa	269.27	Vapor liquid equilibrium data for the hexafluoroethane + carbon dioxide system at temperatures from 253 to 297K and pressures up to 6.5MPa
pvap	2387.00	kPa	283.23	Vapor liquid equilibrium data for the hexafluoroethane + carbon dioxide system at temperatures from 253 to 297K and pressures up to 6.5MPa

pvap	2509.00	kPa	285.23	Vapor liquid equilibrium data for the hexafluoroethane + carbon dioxide system at temperatures from 253 to 297K and pressures up to 6.5MPa
pvap	2636.00	kPa	287.22	Vapor liquid equilibrium data for the hexafluoroethane + carbon dioxide system at temperatures from 253 to 297K and pressures up to 6.5MPa
pvap	2769.00	kPa	289.22	Vapor liquid equilibrium data for the hexafluoroethane + carbon dioxide system at temperatures from 253 to 297K and pressures up to 6.5MPa
pvap	2839.00	kPa	290.23	Vapor liquid equilibrium data for the hexafluoroethane + carbon dioxide system at temperatures from 253 to 297K and pressures up to 6.5MPa
pvap	2873.00	kPa	290.72	Vapor liquid equilibrium data for the hexafluoroethane + carbon dioxide system at temperatures from 253 to 297K and pressures up to 6.5MPa
pvap	2909.00	kPa	291.22	Vapor liquid equilibrium data for the hexafluoroethane + carbon dioxide system at temperatures from 253 to 297K and pressures up to 6.5MPa

pvap	2946.00	kPa	291.72	Vapor liquid equilibrium data for the hexafluoroethane + carbon dioxide system at temperatures from 253 to 297K and pressures up to 6.5MPa
pvap	2982.00	kPa	292.22	Vapor liquid equilibrium data for the hexafluoroethane + carbon dioxide system at temperatures from 253 to 297K and pressures up to 6.5MPa
pvap	1062.90	kPa	253.65	Vapor liquid equilibrium data for the (hexafluoroethane +1,1,1,2-tetrafluoroethane) system at temperatures from 263 to 353K and pressures up to 4.16MPa
pvap	1226.40	kPa	258.46	Vapor liquid equilibrium data for the (hexafluoroethane +1,1,1,2-tetrafluoroethane) system at temperatures from 263 to 353K and pressures up to 4.16MPa
pvap	1415.30	kPa	263.43	Vapor liquid equilibrium data for the (hexafluoroethane +1,1,1,2-tetrafluoroethane) system at temperatures from 263 to 353K and pressures up to 4.16MPa
pvap	1415.30	kPa	263.45	Vapor liquid equilibrium data for the (hexafluoroethane +1,1,1,2-tetrafluoroethane) system at temperatures from 263 to 353K and pressures up to 4.16MPa

pvap	1621.10	kPa	268.33	Vapor liquid equilibrium data for the (hexafluoroethane +1,1,1,2-tetrafluoroethane) system at temperatures from 263 to 353K and pressures up to 4.16MPa
pvap	1850.00	kPa	273.29	Vapor liquid equilibrium data for the (hexafluoroethane +1,1,1,2-tetrafluoroethane) system at temperatures from 263 to 353K and pressures up to 4.16MPa
pvap	1862.00	kPa	273.32	Vapor liquid equilibrium data for the (hexafluoroethane +1,1,1,2-tetrafluoroethane) system at temperatures from 263 to 353K and pressures up to 4.16MPa
pvap	2100.60	kPa	278.19	Vapor liquid equilibrium data for the (hexafluoroethane +1,1,1,2-tetrafluoroethane) system at temperatures from 263 to 353K and pressures up to 4.16MPa
pvap	2392.00	kPa	283.22	Vapor liquid equilibrium data for the (hexafluoroethane +1,1,1,2-tetrafluoroethane) system at temperatures from 263 to 353K and pressures up to 4.16MPa
pvap	2385.80	kPa	283.25	Vapor liquid equilibrium data for the (hexafluoroethane +1,1,1,2-tetrafluoroethane) system at temperatures from 263 to 353K and pressures up to 4.16MPa

pvap	2697.60	kPa	288.20	Vapor liquid equilibrium data for the (hexafluoroethane +1,1,1,2-tetrafluoroethane) system at temperatures from 263 to 353K and pressures up to 4.16MPa
pvap	2697.60	kPa	288.30	Vapor liquid equilibrium data for the (hexafluoroethane +1,1,1,2-tetrafluoroethane) system at temperatures from 263 to 353K and pressures up to 4.16MPa
pvap	1489.00	kPa	265.28	Vapor liquid equilibrium data for the hexafluoroethane + carbon dioxide system at temperatures from 253 to 297K and pressures up to 6.5MPa
pvap	2903.40	kPa	291.22	Vapor liquid equilibrium data for the (hexafluoroethane +1,1,1,2-tetrafluoroethane) system at temperatures from 263 to 353K and pressures up to 4.16MPa
pvap	2976.20	kPa	292.22	Vapor liquid equilibrium data for the (hexafluoroethane +1,1,1,2-tetrafluoroethane) system at temperatures from 263 to 353K and pressures up to 4.16MPa
pvap	1839.00	kPa	273.06	Isothermal vapour-liquid equilibrium data for the binary systems of (CHF3 or C2F6) and n-heptane
pvap	2100.00	kPa	278.07	Isothermal vapour-liquid equilibrium data for the binary systems of (CHF3 or C2F6) and n-heptane

pvap	2388.00	kPa	283.07	Isothermal vapour-liquid equilibrium data for the binary systems of (CHF3 or C2F6) and n-heptane
pvap	2701.00	kPa	288.07	Isothermal vapour-liquid equilibrium data for the binary systems of (CHF3 or C2F6) and n-heptane
pvap	2972.00	kPa	292.07	Isothermal vapour-liquid equilibrium data for the binary systems of (CHF3 or C2F6) and n-heptane
pvap	1240.00	kPa	258.55	Phase Equilibria for Perfluoroethane + (n-Perfluorohexane or n-Perfluorooctane) Binary Systems: Measurement and Modeling
pvap	1300.00	kPa	260.26	Phase Equilibria for Perfluoroethane + (n-Perfluorohexane or n-Perfluorooctane) Binary Systems: Measurement and Modeling
pvap	1390.00	kPa	262.75	Phase Equilibria for Perfluoroethane + (n-Perfluorohexane or n-Perfluorooctane) Binary Systems: Measurement and Modeling
pvap	1480.00	kPa	264.72	Phase Equilibria for Perfluoroethane + (n-Perfluorohexane or n-Perfluorooctane) Binary Systems: Measurement and Modeling

pvap	1550.00	kPa	266.78	Phase Equilibria for Perfluoroethane + (n-Perfluorohexane or n-Perfluorooctane) Binary Systems: Measurement and Modeling
pvap	1650.00	kPa	268.76	Phase Equilibria for Perfluoroethane + (n-Perfluorohexane or n-Perfluorooctane) Binary Systems: Measurement and Modeling
pvap	1730.00	kPa	270.81	Phase Equilibria for Perfluoroethane + (n-Perfluorohexane or n-Perfluorooctane) Binary Systems: Measurement and Modeling
pvap	1830.00	kPa	272.77	Phase Equilibria for Perfluoroethane + (n-Perfluorohexane or n-Perfluorooctane) Binary Systems: Measurement and Modeling
pvap	2000.00	kPa	276.28	Phase Equilibria for Perfluoroethane + (n-Perfluorohexane or n-Perfluorooctane) Binary Systems: Measurement and Modeling
pvap	2160.00	kPa	279.30	Phase Equilibria for Perfluoroethane + (n-Perfluorohexane or n-Perfluorooctane) Binary Systems: Measurement and Modeling

pvap	2330.00	kPa	282.31	Phase Equilibria for Perfluoroethane + (n-Perfluorohexane or n-Perfluorooctane) Binary Systems: Measurement and Modeling
pvap	2470.00	kPa	284.83	Phase Equilibria for Perfluoroethane + (n-Perfluorohexane or n-Perfluorooctane) Binary Systems: Measurement and Modeling
pvap	1851.00	kPa	273.27	Isothermal Vapor-Liquid Equilibrium Data for the Hexafluoroethane (R116) + n-Butane System at Temperatures from 273 to 323 K
pvap	1328.00	kPa	261.29	Vapor liquid equilibrium data for the hexafluoroethane + carbon dioxide system at temperatures from 253 to 297K and pressures up to 6.5MPa
pvap	128.70	kPa	199.64	Vapor-Liquid Equilibrium Measurements for the Ethane + Hexafluoroethane System over a Temperature Range from (199.64 to 242.93) K
pvap	243.40	kPa	213.06	Vapor-Liquid Equilibrium Measurements for the Ethane + Hexafluoroethane System over a Temperature Range from (199.64 to 242.93) K

pvap	453.10	kPa	228.28	Vapor-Liquid Equilibrium Measurements for the Ethane + Hexafluoroethane System over a Temperature Range from (199.64 to 242.93) K
pvap	765.10	kPa	242.93	Vapor-Liquid Equilibrium Measurements for the Ethane + Hexafluoroethane System over a Temperature Range from (199.64 to 242.93) K
pvap	1183.00	kPa	257.31	Vapor liquid equilibrium data for the hexafluoroethane + carbon dioxide system at temperatures from 253 to 297K and pressures up to 6.5MPa
pvap	1048.00	kPa	253.29	Vapor liquid equilibrium data for the hexafluoroethane + carbon dioxide system at temperatures from 253 to 297K and pressures up to 6.5MPa
pvap	928.00	kPa	249.31	Vapor liquid equilibrium data for the hexafluoroethane + carbon dioxide system at temperatures from 253 to 297K and pressures up to 6.5MPa
pvap	816.00	kPa	245.32	Vapor liquid equilibrium data for the hexafluoroethane + carbon dioxide system at temperatures from 253 to 297K and pressures up to 6.5MPa

pvap	769.00	kPa	243.33	Vapor liquid equilibrium data for the hexafluoroethane + carbon dioxide system at temperatures from 253 to 297K and pressures up to 6.5MPa
pvap	1038.20	kPa	252.80	Vapor liquid equilibrium measurements and correlations for an azeotropic system of ethane + hexafluoroethane
pvap	883.00	kPa	247.63	Vapor liquid equilibrium measurements and correlations for an azeotropic system of ethane + hexafluoroethane
pvap	88.50	kPa	192.63	Vapor liquid equilibrium measurements and correlations for an azeotropic system of ethane + hexafluoroethane
pvap	73.90	kPa	189.31	Vapor liquid equilibrium measurements and correlations for an azeotropic system of ethane + hexafluoroethane
pvap	2832.90	kPa	290.22	Vapor liquid equilibrium data for the (hexafluoroethane +1,1,1,2-tetrafluoroethane) system at temperatures from 263 to 353K and pressures up to 4.16MPa
pvap	2708.00	kPa	288.25	Isothermal Vapor-Liquid Equilibrium Data for the Hexafluoroethane (R116) + n-Butane System at Temperatures from 273 to 323 K
rhol	1589.99	kg/m3	195.00	KDB

sfust	35.90	J/mol×K	104.00	NIST Webbook
sfust	15.50	J/mol×K	173.10	NIST Webbook
srf	0.02	N/m	193.20	KDB
svapt	82.88	J/mol×K	194.87	NIST Webbook

Correlations

Information	Value
Property code	pvap
Equation	ln(Pvp) = A + B/(T + C)
Coeff. A	1.44741e+01
Coeff. B	-1.84255e+03
Coeff. C	-8.04800e+00
Temperature range (K), min.	137.93
Temperature range (K), max.	293.03

Information	Value
Property code	pvap
Equation	ln(Pvp) = A + B/T + C*ln(T) + D*T^2
Coeff. A	5.23490e+01
Coeff. B	-3.01585e+03
Coeff. C	-6.23149e+00
Coeff. D	1.57691e-05
Temperature range (K), min.	172.45
Temperature range (K), max.	293.00

Datasets

Speed of sound, m/s

Temperature, K - Gas	Pressure, kPa - Gas	Speed of sound, m/s - Gas
250.80	225.50	124.44
250.80	169.40	125.61
250.80	95.00	127.13
250.80	211.50	124.74

250.80	149.20	126.04
250.80	77.30	127.49
250.80	195.40	125.08
250.80	128.50	126.44
250.80	68.90	127.66
250.80	177.20	125.45
250.80	108.80	126.85
256.20	230.60	125.83
256.20	173.10	126.97
256.20	97.00	128.43
256.20	216.20	126.14
256.20	152.50	127.38
256.20	78.90	128.78
256.20	199.70	126.45
256.20	131.30	127.77
256.20	70.40	128.95
256.20	181.10	126.82
256.20	111.10	128.17
261.50	235.70	127.24
261.50	176.80	128.32
261.50	99.00	129.73
261.50	221.00	127.51
261.50	155.70	128.71
261.50	80.50	130.09
261.50	204.00	127.83
261.50	134.10	129.1
261.50	71.80	130.24
261.50	185.00	128.19
261.50	113.40	129.49
266.80	240.80	128.62
266.80	180.50	129.65
266.80	101.10	131.02
266.80	225.60	128.87
266.80	158.90	130.04
266.80	82.10	131.33
266.80	208.30	129.18
266.80	136.80	130.41
266.80	73.30	131.5
266.80	188.90	129.52
266.80	115.70	130.77
272.10	246.80	129.96
272.10	184.30	131.01
272.10	103.10	132.3
272.10	230.50	130.23

272.10	162.10	131.35
272.10	83.70	132.62
272.10	212.60	130.53
272.10	139.60	131.72
272.10	74.70	132.77
272.10	192.80	130.85
272.10	118.00	132.05
277.40	251.80	131.3
277.40	188.00	132.3
277.40	105.10	133.57
277.40	235.20	131.58
277.40	165.40	132.66
277.40	85.40	133.89
277.40	216.90	131.84
277.40	142.30	132.99
277.40	76.20	134.03
277.40	196.70	132.17
277.40	120.40	133.33
282.70	256.90	132.64
282.70	191.60	133.6
282.70	107.10	134.83
282.70	239.90	132.89
282.70	168.60	133.93
282.70	87.00	135.13
282.70	221.30	133.16
282.70	145.10	134.28
282.70	77.60	135.27
282.70	200.50	133.47
282.70	122.70	134.61
287.90	262.00	133.96
287.90	195.40	134.9
287.90	109.10	136.08
287.90	244.50	134.2
287.90	171.80	135.21
287.90	88.60	136.37
287.90	225.60	134.47
287.90	147.90	135.54
287.90	79.10	136.49
287.90	204.40	134.76
287.90	125.00	135.85
293.20	267.00	135.26
293.20	229.90	135.75
293.20	199.00	136.16
293.20	150.60	136.8

293.20	111.10	137.31
293.20	265.60	135.26
293.20	229.90	135.74
293.20	175.10	136.49
293.20	127.30	137.09
293.20	90.30	137.61
293.20	249.20	135.5
293.20	208.30	136.06
293.20	175.00	136.48
293.20	127.30	137.08
293.20	85.40	137.65
293.20	249.10	135.49
293.20	203.70	136.09
293.20	150.60	136.79
293.20	111.10	137.3
293.20	80.50	137.71
298.50	272.00	136.54
298.50	202.80	137.42
298.50	113.10	138.55
298.50	253.90	136.77
298.50	178.30	137.71
298.50	91.90	138.8
298.50	234.20	137.02
298.50	153.40	138.03
298.50	82.00	138.94
298.50	212.20	137.32
298.50	129.70	138.33
303.70	277.20	137.85
303.70	206.40	138.67
303.70	115.20	139.76
303.70	258.60	138.06
303.70	181.50	138.98
303.70	93.50	140.03
303.70	238.50	138.3
303.70	156.20	139.26
303.70	83.40	140.14
303.70	216.10	138.57
303.70	132.00	139.56
308.90	282.10	139.06
308.90	210.10	139.9
308.90	117.20	140.94
308.90	263.30	139.32
308.90	184.70	140.18
308.90	95.10	141.19

308.90	242.80	139.54
308.90	158.90	140.47
308.90	84.90	141.33
308.90	219.90	139.79
308.90	134.20	140.73
314.10	287.20	140.33
314.10	213.80	141.12
314.10	119.20	142.15
314.10	267.90	140.53
314.10	187.90	141.4
314.10	96.80	142.38
314.10	247.10	140.76
314.10	161.60	141.69
314.10	86.30	142.49
314.10	223.70	140.99
314.10	136.60	141.94
319.30	292.20	141.56
319.30	217.50	142.34
319.30	121.20	143.32
319.30	272.70	141.78
319.30	191.20	142.61
319.30	98.40	143.57
319.30	251.30	141.97
319.30	164.40	142.88
319.30	87.70	143.67
319.30	227.60	142.24
319.30	138.90	143.13
324.50	297.10	142.79
324.50	221.20	143.52
324.50	123.20	144.49
324.50	277.30	142.99
324.50	194.40	143.8
324.50	100.00	144.71
324.50	255.70	143.2
324.50	167.20	144.07
324.50	89.20	144.83
324.50	231.50	143.44
324.50	141.20	144.31

Reference *https://www.doi.org/10.1021/acs.jced.6b00536*

Sources

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Pure Component and Binary Vapor-Liquid Equilibrium + Modeling for Hexafluoropropylene and Perfluoropropane:	https://www.doi.org/10.1021/je900400v
Isothermal Vapor-Liquid Equilibrium Data for the Hexafluoroethane (R116) + Hexafluoropropylene Oxide (HFPO) System at Temperatures from 263 to 323 K:	https://www.doi.org/10.1021/je800849s
Isothermal Bubble Pressure Data for Toluene and Hexafluoroethane for the Binary System of C2F6 and Toluene:	https://www.doi.org/10.1021/acs.jced.8b00825
Phase Equilibria for Perfluoroethane + (n-Perfluorohexane or n-Perfluorooctane) Binary Systems: Measurement and Modeling:	https://www.doi.org/10.1021/acs.jced.6b00409
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
Isothermal Vapor-Liquid Equilibrium Data for the Hexafluoroethane (R116) + n-Butane System at Temperatures from 279 to 323 K:	https://www.doi.org/10.1021/acs.jced.7b00485
Vapor-liquid equilibrium data for the (hexafluoroethane + 1,1,1,2-tetrafluoroethane) system at temperatures from 263 to 353K and pressures up to 4.16MPa:	https://www.doi.org/10.1016/j.fluid.2008.04.002
Experimental determination of the critical loci for R-23 + propane or R-125 + propane mixtures:	https://www.doi.org/10.1016/j.jct.2016.12.021
Speed of Sound Data in Pure Refrigerants R-116 and R-218 and Their Mixtures: Experiment and Modeling:	https://www.doi.org/10.1021/acs.jced.6b00536
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C76164&Units=SI
Vapor-Liquid Equilibrium Measurements for the Ethane + Hexafluoroethane System over a Temperature Range from (199.64 to 242.93) K:	https://www.doi.org/10.1021/je050281w
KDB:	https://www.cheric.org/files/research/kdb/mol/mol1547.mol
PVTx Measurements for the R116 + CO2 and R41 + CO2 Systems. New Isochoric Apparatus:	https://www.doi.org/10.1021/je049939g
Isothermal Vapor-Liquid Equilibrium Data for Binary Mixtures of Hexafluoroethane (R116) + n-Pentane at Two Temperatures, 288 and 298 K:	https://www.doi.org/10.1021/acs.jced.7b00836
Vapor-liquid equilibrium measurements and correlations for an azeotropic system of hexafluoroethane + n-hexane:	https://www.doi.org/10.1016/j.fluid.2005.12.004
Isothermal vapour-liquid equilibrium data for the binary systems of (CHF3 or C2F6) and n-heptane:	https://www.doi.org/10.1016/j.jct.2016.06.031
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Velocity of sound in perfluoropropane (C3F8), perfluoroethane (C2F6) and their mixtures:	https://www.doi.org/10.1016/j.fluid.2012.10.002
An isothermal vapour-liquid equilibrium data for the binary systems of CHF3 with (n-nonane, n-decane, or n-undecane) and C2F6 with (n-nonane or n-decane):	https://www.doi.org/10.1016/j.fluid.2018.02.015

Legend

cpg:	Ideal gas heat capacity
cpl:	Liquid phase heat capacity
dm:	Dipole Moment
gf:	Standard Gibbs free energy of formation
hf:	Enthalpy of formation at standard conditions
hfus:	Enthalpy of fusion at standard conditions
hfust:	Enthalpy of fusion at a given temperature
hsubt:	Enthalpy of sublimation at a given temperature

hvap:	Enthalpy of vaporization at standard conditions
hvapt:	Enthalpy of vaporization at a given temperature
ie:	Ionization energy
log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
pc:	Critical Pressure
pt:	Triple Point Pressure
pvap:	Vapor pressure
rhoc:	Critical density
rhol:	Liquid Density
rinpol:	Non-polar retention indices
sfust:	Entropy of fusion at a given temperature
sl:	Liquid phase molar entropy at standard conditions
speedsl:	Speed of sound in fluid
srf:	Surface Tension
svapt:	Entropy of vaporization at a given temperature
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
tt:	Triple Point Temperature
vc:	Critical Volume
zc:	Critical Compressibility